## Amendments to the Claims

The following listing of claims will replace all prior versions, and listings, of claims in the application:

- 1-9. (Canceled).
- 10. (Previously presented) A compound having the formula:

$$\begin{array}{c|c}
 & O & O \\
 & N & \\
 & R^1 & \\
 & H & \\
\end{array}$$

wherein:

 $R^1$  is H,  $(C_1-C_8)$ alkyl,  $(C_3-C_7)$ cycloalkyl,  $(C_2-C_8)$ alkenyl,  $(C_2-C_8)$ alkynyl, benzyl, aryl,  $(C_0-C_4)$ alkyl– $(C_1-C_6)$ heterocycloalkyl,  $(C_0-C_4)$ alkyl– $(C_2-C_5)$ heteroaryl,  $C(O)R^3$ ,  $C(S)R^3$ ,  $C(O)OR^4$ ,  $(C_1-C_8)$ alkyl– $N(R^6)_2$ ,  $(C_1-C_8)$ alkyl– $OR^5$ ,  $(C_1-C_8)$ alkyl– $O(CO)R^5$ ,  $C(O)NHR^3$ ,  $C(S)NHR^3$ ,  $C(O)NR^3R^3$ ,  $C(S)NR^3R^3$  or  $(C_1-C_8)$ alkyl– $O(CO)R^5$ ;

 $R^2$  is H or  $(C_1-C_8)$ alkyl;

 $R^3$  and  $R^3$ ' are independently  $(C_1-C_8)$ alkyl,  $(C_3-C_7)$ cycloalkyl,  $(C_2-C_8)$ alkenyl,  $(C_2-C_8)$ alkynyl, benzyl, aryl,  $(C_0-C_4)$ alkyl- $(C_1-C_6)$ heterocycloalkyl,  $(C_0-C_4)$ alkyl- $(C_2-C_5)$ heteroaryl,  $(C_0-C_8)$ alkyl- $N(R^6)_2$ ,  $(C_1-C_8)$ alkyl- $O(C^5)$ ,  $(C_1-C_8)$ alkyl- $O(C^5)$ , or  $O(C^5)$ ;

 $R^4 \text{ is } (C_1-C_8) \text{alkyl, } (C_2-C_8) \text{alkenyl, } (C_2-C_8) \text{alkynyl, } (C_1-C_4) \text{alkyl-OR}^5, \text{ benzyl, aryl, } (C_0-C_4) \text{alkyl-} (C_1-C_6) \text{heterocycloalkyl, or } (C_0-C_4) \text{alkyl-} (C_2-C_5) \text{heteroaryl; } \\ R^5 \text{ is } (C_1-C_8) \text{alkyl, } (C_2-C_8) \text{alkenyl, } (C_2-C_8) \text{alkynyl, benzyl, aryl, or } (C_2-C_5) \text{heteroaryl; } \\ C_5) \text{heteroaryl; } C_5 \text{ is } (C_1-C_8) \text{alkynyl, } (C_2-C_8) \text{alkynyl, benzyl, aryl, or } (C_2-C_8) \text{alkynyl, } \\ C_5) \text{heteroaryl; } C_5 \text{ is } (C_1-C_8) \text{ alkynyl, } (C_2-C_8) \text{ alkynyl, } (C_2-C_8) \text{ alkynyl, } \\ C_5) \text{ is } (C_1-C_8) \text{ alkyl, } (C_2-C_8) \text{ alkynyl, } (C_2-C_8) \text{ alkynyl, } (C_2-C_8) \text{ alkynyl, } (C_2-C_8) \text{ alkynyl, } \\ C_5) \text{ is } (C_1-C_8) \text{ alkyl, } (C_2-C_8) \text{ alkynyl, } (C_2-C_8) \text{ alkynyl, } (C_2-C_8) \text{ alkynyl, } \\ C_5) \text{ is } (C_1-C_8) \text{ alkyl, } (C_2-C_8) \text{ alkynyl, } (C_2-C_8) \text{ alkynyl, } (C_2-C_8) \text{ alkynyl, } \\ C_5) \text{ is } (C_1-C_8) \text{ alkyl, } (C_2-C_8) \text{ alkynyl, } \\ C_5) \text{ is } (C_1-C_8) \text{ alkyl, } (C_2-C_8) \text{ alkynyl, } (C_2-C_8) \text$ 

each occurrence of  $R^6$  is independently H,  $(C_1-C_8)$ alkyl,  $(C_2-C_8)$ alkenyl,  $(C_2-C_8)$ alkynyl, benzyl, aryl,  $(C_2-C_5)$ heteroaryl, or  $(C_0-C_8)$ alkyl- $C(O)O-R^5$  or the  $R^6$  groups can join to form a heterocycloalkyl group; and the \* represents a chiral-carbon center.

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11. (Original) A compound of claim 10, wherein R<sup>1</sup> is H, (C<sub>1</sub>-C<sub>4</sub>)alkyl, CH<sub>2</sub>OCH<sub>3</sub>, CH<sub>2</sub>CH<sub>2</sub>OCH<sub>3</sub>, or

$$\sim$$
 CH<sub>2</sub>  $\sim$  Or  $\sim$  CH<sub>2</sub>  $\sim$  R<sup>7</sup>,

wherein Q is O or S, and each occurrence of  $R^7$  is independently H,(C<sub>1</sub>–C<sub>8</sub>)alkyl, (C<sub>3</sub>–C<sub>7</sub>)cycloalkyl, (C<sub>2</sub>–C<sub>8</sub>)alkenyl, (C<sub>2</sub>–C<sub>8</sub>)alkynyl, benzyl, aryl, halogen, (C<sub>0</sub>–C<sub>4</sub>)alkyl–(C<sub>1</sub>–C<sub>6</sub>)heterocycloalkyl, (C<sub>0</sub>–C<sub>4</sub>)alkyl–(C<sub>2</sub>–C<sub>5</sub>)heteroaryl, (C<sub>0</sub>–C<sub>8</sub>)alkyl–N( $R^6$ )<sub>2</sub>, (C<sub>1</sub>–C<sub>8</sub>)alkyl–OR<sup>5</sup>, (C<sub>1</sub>–C<sub>8</sub>)alkyl–O(O)OR<sup>5</sup>, or C(O)OR<sup>5</sup>, or adjacent occurrences of  $R^7$  can be taken together to form a bicyclic alkyl or aryl ring.

- 12. (Original) A compound of claim 10, wherein  $R^1$  is  $C(O)R^3$ .
- 13. (Original) A compound of claim 10, wherein R<sup>1</sup> is C(O)OR<sup>4</sup>.
- 14. (Previously presented) A compound having the formula:

wherein:

 $R^1$  is H,  $(C_1-C_8)$ alkyl,  $(C_3-C_7)$ cycloalkyl,  $(C_2-C_8)$ alkenyl,  $(C_2-C_8)$ alkynyl, benzyl, aryl,  $(C_0-C_4)$ alkyl- $(C_1-C_6)$ heterocycloalkyl,  $(C_0-C_4)$ alkyl- $(C_2-C_5)$ heteroaryl,  $C(O)R^3$ ,  $C(S)R^3$ ,  $C(O)OR^4$ ,  $(C_1-C_8)$ alkyl- $N(R^6)_2$ ,  $(C_1-C_8)$ alkyl- $OR^5$ ,  $(C_1-C_8)$ alkyl- $C(O)OR^5$ ,  $C(O)NHR^3$ ,  $C(S)NHR^3$ ,  $C(O)NR^3R^3$ ,  $C(S)NR^3R^3$  or  $(C_1-C_8)$ alkyl- $O(CO)R^5$ ;

 $R^2$  is H or  $(C_1-C_8)$ alkyl;

 $R^3$  and  $R^{3'}$  are independently  $(C_1-C_8)$ alkyl,  $(C_3-C_7)$ cycloalkyl,  $(C_2-C_8)$ alkenyl,  $(C_2-C_8)$ alkynyl, benzyl, aryl,  $(C_0-C_4)$ alkyl- $(C_1-C_6)$ heterocycloalkyl,  $(C_0-C_4)$ alkyl- $(C_2-C_5)$ heteroaryl,  $(C_0-C_8)$ alkyl- $(C_1-C_8)$ alkyl- $(C_1-C_$ 

 $R^4$  is  $(C_1-C_8)$ alkyl,  $(C_2-C_8)$ alkenyl,  $(C_2-C_8)$ alkynyl,  $(C_1-C_4)$ alkyl- $(C_1-C_6)$ heterocycloalkyl, or  $(C_0-C_4)$ alkyl- $(C_2-C_5)$ heteroaryl;  $R^5$  is  $(C_1-C_8)$ alkyl,  $(C_2-C_8)$ alkenyl,  $(C_2-C_8)$ alkynyl, benzyl, aryl, or  $(C_2-C_5)$ heteroaryl;

each occurrence of  $R^6$  is independently H,  $(C_1-C_8)$ alkyl,  $(C_2-C_8)$ alkenyl,  $(C_2-C_8)$ alkynyl, benzyl, aryl,  $(C_2-C_5)$ heteroaryl, or  $(C_0-C_8)$ alkyl- $C(O)O-R^5$  or the  $R^6$  groups can join to form a heterocycloalkyl group; and the \* represents a chiral-carbon center.

15. (Original) A compound of claim 14, wherein R<sup>1</sup> is H, (C<sub>1</sub>-C<sub>4</sub>)alkyl, CH<sub>2</sub>OCH<sub>3</sub>, CH<sub>2</sub>CH<sub>2</sub>OCH<sub>3</sub>, or

wherein Q is O or S, and each occurrence of  $R^7$  is independently H,(C<sub>1</sub>–C<sub>8</sub>)alkyl, (C<sub>3</sub>–C<sub>7</sub>)cycloalkyl, (C<sub>2</sub>–C<sub>8</sub>)alkenyl, (C<sub>2</sub>–C<sub>8</sub>)alkynyl, benzyl, aryl, halogen, (C<sub>0</sub>–C<sub>4</sub>)alkyl–(C<sub>1</sub>–C<sub>6</sub>)heterocycloalkyl, (C<sub>0</sub>–C<sub>4</sub>)alkyl–(C<sub>2</sub>–C<sub>5</sub>)heteroaryl, (C<sub>0</sub>–C<sub>8</sub>)alkyl–N( $R^6$ )<sub>2</sub>, (C<sub>1</sub>–C<sub>8</sub>)alkyl–OR<sup>5</sup>, (C<sub>1</sub>–C<sub>8</sub>)alkyl–O(CO)R<sup>5</sup>, or C(O)OR<sup>5</sup>, or adjacent occurrences of  $R^7$  can be taken together to form a bicyclic alkyl or aryl ring.

- 16. (Original) A compound of claim 14, wherein  $R^1$  is  $C(O)R^3$ .
- 17. (Original) A compound of claim 14, wherein R<sup>1</sup> is C(O)OR<sup>4</sup>.
- 18. (Previously presented) A compound having the formula:

wherein:

 $R^1$  is H,  $(C_1-C_8)$ alkyl,  $(C_3-C_7)$ cycloalkyl,  $(C_2-C_8)$ alkenyl,  $(C_2-C_8)$ alkynyl, benzyl, aryl,  $(C_0-C_4)$ alkyl- $(C_1-C_6)$ heterocycloalkyl,  $(C_0-C_4)$ alkyl- $(C_2-C_5)$ heteroaryl,  $C(O)R^3$ ,  $C(S)R^3$ ,  $C(O)OR^4$ ,  $(C_1-C_8)$ alkyl- $N(R^6)_2$ ,  $(C_1-C_8)$ alkyl- $OR^5$ ,  $(C_1-C_8)$ alkyl- $C(O)OR^5$ ,  $C(O)NHR^3$ ,  $C(S)NHR^3$ ,  $C(O)NR^3R^3$ ,  $C(S)NR^3R^3$  or  $(C_1-C_8)$ alkyl- $O(CO)R^5$ :

 $R^2$  is H or  $(C_1-C_8)$  alkyl;

 $R^3$  and  $R^{3'}$  are independently (C<sub>1</sub>–C<sub>8</sub>)alkyl, (C<sub>3</sub>–C<sub>7</sub>)cycloalkyl, (C<sub>2</sub>–C<sub>8</sub>)alkenyl, (C<sub>2</sub>–C<sub>8</sub>)alkynyl, benzyl, aryl, (C<sub>0</sub>–C<sub>4</sub>)alkyl–(C<sub>1</sub>–C<sub>6</sub>)heterocycloalkyl, (C<sub>0</sub>–C<sub>4</sub>)alkyl–(C<sub>2</sub>–C<sub>5</sub>)heteroaryl, (C<sub>0</sub>–C<sub>8</sub>)alkyl–N( $R^6$ )<sub>2</sub>, (C<sub>1</sub>–C<sub>8</sub>)alkyl–OR<sup>5</sup>, (C<sub>1</sub>–C<sub>8</sub>)alkyl–C(O)OR<sup>5</sup>, (C<sub>1</sub>–C<sub>8</sub>)alkyl–O(CO)R<sup>5</sup>, or C(O)OR<sup>5</sup>;

 $R^4 \text{ is } (C_1-C_8)\text{alkyl}, (C_2-C_8)\text{alkenyl}, (C_2-C_8)\text{alkynyl}, (C_1-C_4)\text{alkyl}-OR^5, \text{ benzyl}, \text{ aryl}, \\ (C_0-C_4)\text{alkyl}-(C_1-C_6)\text{heterocycloalkyl}, \text{ or } (C_0-C_4)\text{alkyl}-(C_2-C_5)\text{heteroaryl}; \\ R^5 \text{ is } (C_1-C_8)\text{alkyl}, (C_2-C_8)\text{alkenyl}, (C_2-C_8)\text{alkynyl}, \text{ benzyl}, \text{ aryl}, \text{ or } (C_2-C_5)\text{heteroaryl}; \\ C_5)\text{heteroaryl};$ 

each occurrence of  $R^6$  is independently H,  $(C_1-C_8)$ alkyl,  $(C_2-C_8)$ alkenyl,  $(C_2-C_8)$ alkynyl, benzyl, aryl,  $(C_2-C_5)$ heteroaryl, or  $(C_0-C_8)$ alkyl- $C(O)O-R^5$  or the  $R^6$  groups can join to form a heterocycloalkyl group; and the \* represents a chiral-carbon center.

19. (Original) A compound of claim 18, wherein R<sup>1</sup> is H, (C<sub>1</sub>-C<sub>4</sub>)alkyl, CH<sub>2</sub>OCH<sub>3</sub>, CH<sub>2</sub>CCH<sub>2</sub>OCH<sub>3</sub> or

$$\text{or } \text{vcH}_2 \longrightarrow \text{or } \text{vcH}_2 \longrightarrow \text{R}^7$$

wherein Q is O or S, and each occurrence of  $R^7$  is independently H,(C<sub>1</sub>–C<sub>8</sub>)alkyl, (C<sub>3</sub>–C<sub>7</sub>)cycloalkyl, (C<sub>2</sub>–C<sub>8</sub>)alkenyl, (C<sub>2</sub>–C<sub>8</sub>)alkynyl, benzyl, aryl, halogen, (C<sub>0</sub>–C<sub>4</sub>)alkyl–(C<sub>1</sub>–C<sub>6</sub>)heterocycloalkyl, (C<sub>0</sub>–C<sub>4</sub>)alkyl–(C<sub>2</sub>–C<sub>5</sub>)heteroaryl, (C<sub>0</sub>–C<sub>8</sub>)alkyl–N( $R^6$ )<sub>2</sub>, (C<sub>1</sub>–C<sub>8</sub>)alkyl–OR<sup>5</sup>, (C<sub>1</sub>–C<sub>8</sub>)alkyl–O(CO)R<sup>5</sup>, or C(O)OR<sup>5</sup>, or adjacent occurrences of  $R^7$  can be taken together to form a bicyclic alkyl or aryl ring.

20. (Original) A compound of claim 18, wherein  $R^1$  is  $C(O)R^3$ .

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21. (Original) A compound of claim 18, wherein R<sup>1</sup> is C(O)OR<sup>4</sup>.

22-100. (Canceled).

(Currently amended) A compound of claim 10, which is: N-[2-(2,6-101. dioxo-piperidin 3-yl) 1,3 dioxo 2,3 duhydro 1H isoindol 4-yl-methyl] acetamide; N {[2-(2,6-dioxo(3-piperidyl))-1,3-dioxoisoindolin-4-yl]methyl}cyclopropylcarboxamide; 1-tert-butyl-3-[2-(2,6-dioxo-piperidin-3-yl)-1,3-dioxo-2,3-dihydro-1Hisoindol 4-ylmethyl] urea; N-{[2 (2,6-dioxo(3-piperidyl))-1,3-dioxoisoindolin-4yl|methyl}-3,3-dimethylbutanamide; N-{[2-(2,6-dioxo(3-piperidyl))-1,3dioxoisoindolin 4 yl]methyl} propanamide; N-{[2 (2,6 dioxo(3 piperidyl)) 1,3dioxoisoindolin 4-yl]methyl}-3-pyridylcarboxamide; N-{[2-(2,6-dioxo(3-piperidyl))-1.3-dioxoisoindolin 4-yl]methyl}heptanamide; N {[2 (2,6 dioxo(3 piperidyl)) 1,3dioxoisoindolin 4-yl]methyl}-2-furylcarboxamide; 2-amino N-{[2-(2,6-dioxo(3piperidyl)) 1,3 dioxoisoindolin 4 yl]methyl}acetamide; N {[2 (2,6 dioxo(3piperidyl)) 1,3-dioxoisoindolin 4 yl]methyl} 2 thienylcarboxamide; N {[2 (2,6dioxo(3-piperidyl)) 1,3-dioxoindolin 4yl]methyl] (ethylamino)carboxamide; N-{[2-(2,6-dioxo(3-piperidyl)) 1,3-dioxoisoindolin 4-yl]methyl}butanamide; N {[2-(2,6-dioxo(3-piperidyl)) 1,3-dioxoisoindolin 4-yl]methyl}butanamide; N {[2-(2,6-dioxo(3-piperidyl)] 1,3-dioxoisoindolin 4-yl]methyll 1,3-dioxoisoindolin 4-yl]methyll 1,3-dioxoisoindolin 4-yl]methyll 1,3-dioxoisoindolin 4-yl]methyll 1,3-dioxoisoin dioxo(3-piperidyl)) 1,3-dioxoisoindolin 4-yl]methyl} 2-pyridylcarboxamide; N-{[2-(2,6-dioxo(3-piperidyl)) 1,3-dioxoisoindolin 4-yl]methyl}undecamide; N-{[2-(2,6-dioxo(3-piperidyl)] 1,3-dioxoisoindolin 4-yl]methyllundecamide; N-{[2-(2,6-dioxo(3-piperidyl)] 1,3-dioxoisoindolin 4-yl]methyllundecamide; N-{[2-(2,6-dioxo(3-piperidyl)] 1,3-dioxoisoindolin 4-yl]methyllundecamide; N-{[2-(2,6-dioxo(3-piperidy dioxo(3-piperidyl)) 1,3-dioxoisoindolin-4-yl]methyl}2 methylpropanamide; N-{[2-(2.6 dioxo(3 piperidyl)) 1.3 dioxoisoindolin 4 yl]methyl]cyclopentylcarboxamide; N-{[2 (2,6-dioxo(3-piperidyl)) 1,3-dioxoisoindolin-4-yl]methyl} eyclohexylcarboxamide; N {[2 (2,6 dioxo(3 piperidyl)) 1,3 dioxoisoindolin-4yl]methyl}(butylamino)carboxamide; N-{[2-(2,6-dioxo(3-piperidyl))-1,3dioxoisoindolin 4 yl]methyl}(propylamino)carboxamide; N {[2-(2,6 dioxo(3piperidyl))-1,3 dioxoisoindolin 4 yl]methyl}[(methylethylamino)] carboxamide; N-{[2-(2,6-dioxo(3-piperidyl))-1,3-dioxoisoindolin-4yl]methyl}(octylamino)carboxamide; N-{[2-(2,6 dioxo(3-piperidyl)) 1,3dioxoisoindolin-4-yl]methyl}(cyclopropylamino)carboxamide; or N-{[2-(2,6dioxo(3-piperidyl)) 1,3-dioxoisoindolin 4-yl]methyl}(diethylamino)carboxamide.

(Claim 101 continued on next page)

H 3C N O O N H O	0 0 N H O	NH NHO
H <sub>3</sub> C O N H O N H O	H <sub>3</sub> C NHO	
0 0 N H O	N — N H O	H N H 2
0 0 N H O	0 0 N H O O N H O	H N O O N H O
0 0 N H O	H N O N H	N N H O
0 0 N H O	0 0 N H O	O O N H O O N H O O N H O O N H O O O O
	N N N N N N N N N N N N N N N N N N N	or O

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102. (Currently amended) A compound of claim 10, which is:  $\frac{2-(2,6)}{(2,6)}$ dioxo-piperidin-3-yl) 1,3-dioxo-2,3-dihydro-1H-isoindol-4-yl-methyl]-carbamic acid tert-butyl ester; 4-(aminomethyl)-2-(2,6-dioxo(3-Piperidyl)) isoindoline-1,3-dione; [2-(2,6-dioxo-piperidin 3-yl) 1,3-dioxo-2,3-dihydro-1H-isoindol-4-yl-methyl] carbamic acid ethyl ester; [2 (2,6-dioxo-piperidin 3-yl)-1,3-dioxo-2,3-dihydro-1Hisoindol 4 yl methyl] carbamic acid benzyl ester; 2 (dimethylamino) N {[2 (2,6dioxo(3-piperidyl)) 1,3-dioxoisoindolin 4-yl]methyl}acetamide; ethyl 6-(3N-{[2-(2,6dioxo(3-piperidyl)) 1,3-dioxoisoindolin 4-yl]methyl]carbamoyl)hexanoate; 3-[(tertbutoxy)carbonylaminol-N-{[2-(2,6-dioxo(3-piperidyl))-1,3-dioxoisoindolin-4yl]methyl}propanamide; 3-amino-N-{[2-(2,6-dioxo(3-piperidyl)) 1,3dioxoisoindolin 4 yl]methyl}propanamide; N-{[2-(2,6-dioxo(3-piperidyl))-1,3dioxoisoindolin 4 yl]methyl} 2 methoxyacetamide; (N {[2 (2,6 dioxo(3 piperidyl)) 1,3-dioxoisoindolin-4-yl]methyl}carbamoyl)methyl acetate; ethyl 2-[N-{[2 (2,6dioxo(3-piperidyl)) 1,3-dioxoisoindolin 4-yl]methyl]carbamoyl)amino]acetate; 7amino-N-{[2-(2,6-dioxo(3-piperidyl)) 1,3-dioxoisoindolin 4-yl]methyl}heptanamide; N-{[2-(2,6-dioxo(3-piperidyl)) 1,3-dioxoisoindolin 4-yl]methyl}benzamide; N-{[2-(2,6 dioxo(3-piperidyl))-1,3 dioxoisoindolin 4-yl]methyl}phenylacetamide; N-{[2-(2,6-dioxo(3-piperidyl))-1,3-dioxoisoindolin-4yl]methyl}(phenylamino)carboxamide; N-{[2-(2,6-dioxo(3-piperidyl))-1,3dioxoisoindolin 4 yl]methyl}(benzylamino)carboxamide, 2 (2,6 dioxo piperidin 3yl) 4-{[(furan-2-ylmethyl)-amino-methyl}-isoindole-1,3-dione; N-{[2-(2,6-dioxo(3piperidyl)) 1,3 dioxo 2,3 dihydro 1H-isoindol 4 ylmethyl] isonicotinamide; 2 (2,6dioxo(3-piperidyl))-4-({[(cyclohexylamino)thioxomethyl]amino}methyl)isoindole-1,3-dione; 2-(2,6-dioxo(3-piperidyl))-4-({[(ethylamino)thioxomethyl]amino}methyl)isoindole-1,3-dione; 2-(2,6-dioxo(3piperidyl))-4-({[(propylamino)thioxomethyl]amino} methyl)isoindole-1,3-dione; N-{[2-(2,6-dioxo(3-piperidyl))-1,3-dioxoisoindolin-4yl]methyl}(cyclopentylamino)carboxamide; N-{[2-(2,6-dioxo(3-piperidyl))-1,3-

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dioxoisoindolin 4-yl]methyl}(3-pyridylamino)carboxamide; N-{[2-(2,6-dioxo(3-piperidyl))-1,3-dioxoisoindolin 4-yl]methyl}piperidylcarboxamide; or piperazine-1-carboxylic acid [2-(2,6-dioxo-piperidin-3-yl)-1,3-dioxo-2,3-dihydro-1H-isoindol-4-ylmethyl]-amide.

t-B u 0 N H 0	0 0 N H O	H <sub>3</sub> C 0 H 0 0 N H 0
	H <sub>3</sub> C 0 N H O	0 0 N H O O C H 3
	H N O N H 2	0 C H 3
0 0 N H O C H 3	0 0 0 H 0 0 C H 3	H N O N H 2
0 0 N H O	H N O O N H O	
N N H O	N N N N N N N N N N N N N N N N N N N	

H N H O	S N N N N N N N N N N N N N N N N N N N	
S H	H N N O N H O	N H O N H O
0 0 N H O	or O O O O O O O O O O O O O O O O O O O	0 0 N H O N H O N H

103. (Currently amended) A compound of claim 14, which is: N-{2-(2,6-dioxo-piperidin-3-yl) 1-oxo-2,3-dihydro-1H-isoindol-4-ylmethyl} acetamide; N-{[2-(2,6-dioxo(3-piperidyl)) 1-oxoisoindolin-4-yl]methyl}cyclopropylcarboxamide; or N-{[2-(2,6-dioxo(3-piperidyl)) 1-oxoisoindolin-4-yl]methyl}(ethylamino)carboxamide.

104-105. (Canceled)